

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042319\
 Data File : BF113937.D
 Acq On : 23 Apr 2019 22:45
 Operator : JU/SJ
 Sample : K2515-19RE
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-13RE

Manual Integrations
 APPROVED

Sohil
 4/24/2019 11:51:33 AM

Quant Time: Apr 24 07:38:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	94061	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	300143	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	127754	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	233348	20.00	ng	0.00
76) Chrysene-d12	14.05	240	268117	20.00	ng	-0.02
87) Perylene-d12	15.53	264	230333	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1045292	190.07	ng	0.01
7) Phenol-d6	6.53	99	1308589	189.66	ng	0.00
23) Nitrobenzene-d5	7.46	82	758099	155.95	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	312321	200.68	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1336065	163.57	ng	0.00
79) Terphenyl-d14	13.00	244	1550998	118.60	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.64	163	99036	10.719	ng	99
75) Fluoranthene	12.63	202	57514	4.495	ng	95
78) Pyrene	12.86	202	53131	2.834	ng	97
81) Benzo(a)anthracene	14.04	228	34760	2.200	ng	95
83) Chrysene	14.07	228	35616	2.315	ng	99
88) Benzo(b)fluoranthene	15.10	252	45575m	3.127	ng	
90) Benzo(a)pyrene	15.47	252	29062	2.306	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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